Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079422.D
Acq On : 22 May 2015 2:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 22 05:02:09 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 22 05:01:56 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	58	-0.06
2		1,4-Dioxane	40.000	49.462	-23.7	73	-0.08
3		Pyridine	40.000	41.285	-3.2	64	-0.09
4		n-Nitrosodimethylamine	40.000	37.318	6.7	56	-0.08
5	S	2-Fluorophenol	80.000	85.183	-6.5	64	-0.07
6		Aniline	40.000	40.680	-1.7	61	-0.06
7	S	Phenol-d6	80.000	83.497	-4.4	65	-0.06
8		2-Chlorophenol	40.000	43.579	-8.9	69	-0.07
9		Benzaldehyde	40.000	37.580	6.1	57	-0.06
10	C	Phenol	40.000	42.665	-6.7	64	-0.06
11		bis(2-Chloroethyl)ether	40.000	42.308	-5.8	67	-0.06
12		1,3-Dichlorobenzene	40.000	42.625	-6.6	67	-0.06
13	C	1,4-Dichlorobenzene	40.000	42.371	-5.9	66	-0.06
14		1,2-Dichlorobenzene	40.000	42.914	-7.3	66	-0.06
15		Benzyl Alcohol	40.000	38.074	4.8	59	-0.06
16		2,2'-oxybis(1-Chloropropane	40.000	40.048	-0.1	63	-0.06
17		2-Methylphenol	40.000	43.262	-8.2	67	-0.06
18		Hexachloroethane	40.000	42.897	-7.2	64	-0.06
19	P	n-Nitroso-di-n-propylamine	40.000	41.304	-3.3	61	-0.06
20		3+4-Methylphenols	40.000	42.673	-6.7	64	-0.06
21	I	Naphthalene-d8	20.000	20.000	0.0	59	-0.06
22		Acetophenone	40.000	42.016	-5.0	67	-0.06
23	S	Nitrobenzene-d5	80.000	84.471	-5.6	65	-0.06
24		Nitrobenzene	40.000	42.193	-5.5	68	-0.06
25		Isophorone	40.000	41.896	-4.7	64	-0.06
26	C	2-Nitrophenol	40.000	40.776	-1.9	60	-0.06
27		2,4-Dimethylphenol	40.000	40.319	-0.8	61	-0.05
28		bis(2-Chloroethoxy)methane	40.000	42.731	-6.8	64	-0.06
29	C	2,4-Dichlorophenol	40.000	43.466	-8.7	67	-0.06
30		1,2,4-Trichlorobenzene	40.000	43.096	-7.7	67	-0.06
31		Naphthalene	40.000	43.875	-9.7	69	-0.06
32		Benzoic acid	40.000	41.196	-3.0	62	-0.07
33		4-Chloroaniline	40.000	43.174	-7.9	68	-0.06
34	C	Hexachlorobutadiene	40.000	42.027	-5.1	67	-0.06
35		Caprolactam	40.000	42.188	-5.5	64	-0.07
36	C	4-Chloro-3-methylphenol	40.000	44.911	-12.3	65	-0.06
37		2-Methylnaphthalene	40.000	43.442	-8.6	67	-0.06
38	I	Acenaphthene-d10	20.000	20.000	0.0	58	-0.07
39		1,2,4,5-Tetrachlorobenzene	40.000	42.470	-6.2	67	-0.07
40	P	Hexachlorocyclopentadiene	40.000	45.064	-12.7	68	-0.06
41	S	2,4,6-Tribromophenol	80.000	86.304	-7.9	63	-0.07
42	C	2,4,6-Trichlorophenol	40.000	41.729	-4.3	62	-0.07
43		2,4,5-Trichlorophenol	40.000	41.017	-2.5	62	-0.06
44	S	2-Fluorobiphenyl	80.000	88.190	-10.2	67	-0.06
45		1,1'-Biphenyl	40.000	42.889	-7.2	67	-0.06
46		2-Chloronaphthalene	40.000	43.451	-8.6	69	-0.06
47		2-Nitroaniline	40.000	41.775	-4.4	62	-0.06
48		Acenaphthylene	40.000	44.459	-11.1	69	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079422.D
Acq On : 22 May 2015 2:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 22 05:02:09 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 22 05:01:56 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
49		Dimethylphthalate	40.000	41.466	-3.7	66	-0.07
50		2,6-Dinitrotoluene	40.000	44.273	-10.7	67	-0.06
51	C	Acenaphthene	40.000	42.051	-5.1	64	-0.06
52		3-Nitroaniline	40.000	43.774	-9.4	67	-0.07
53	P	2,4-Dinitrophenol	40.000	26.846	32.9#	32	-0.06
54		Dibenzofuran	40.000	43.023	-7.6	66	-0.06
55	P	4-Nitrophenol	40.000	29.109	27.2#	45	-0.05
56		2,4-Dinitrotoluene	40.000	43.132	-7.8	62	-0.07
57		Fluorene	40.000	43.574	-8.9	69	-0.06
58		2,3,4,6-Tetrachlorophenol	40.000	38.559	3.6	57	-0.06
59		Diethylphthalate	40.000	43.029	-7.6	67	-0.06
60		4-Chlorophenyl-phenylether	40.000	42.655	-6.6	66	-0.06
61		4-Nitroaniline	40.000	44.724	-11.8	66	-0.07
62		Azobenzene	40.000	42.103	-5.3	63	-0.06
63	I	Phenanthrene-d10	20.000	20.000	0.0	59	-0.07
64		4,6-Dinitro-2-methylphenol	40.000	33.541	16.1	47	-0.07
65	c	n-Nitrosodiphenylamine	40.000	42.608	-6.5	65	-0.06
66		4-Bromophenyl-phenylether	40.000	42.042	-5.1	67	-0.06
67		Hexachlorobenzene	40.000	41.617	-4.0	67	-0.07
68		Atrazine	40.000	39.930	0.2	63	-0.07
69	C	Pentachlorophenol	40.000	30.443	23.9#	45	-0.07
70		Phenanthrene	40.000	41.966	-4.9	65	-0.06
71		Anthracene	40.000	41.954	-4.9	66	-0.07
72		Carbazole	40.000	42.939	-7.3	67	-0.06
73		Di-n-butylphthalate	40.000	41.732	-4.3	62	-0.07
74	C	Fluoranthene	40.000	42.821	-7.1	66	-0.07
75	I	Chrysene-d12	20.000	20.000	0.0	55	-0.09
76		Benzidine	40.000	43.215	-8.0	58	-0.07
77		Pyrene	40.000	44.978	-12.4	66	-0.07
78	S	Terphenyl-d14	80.000	89.329	-11.7	65	-0.07
79		Butylbenzylphthalate	40.000	45.192	-13.0	61	-0.07
80		Benzo(a)anthracene	40.000	43.037	-7.6	62	-0.09
81		3,3'-Dichlorobenzidine	40.000	43.292	-8.2	60	-0.08
82		Chrysene	40.000	43.405	-8.5	65	-0.09
83		Bis(2-ethylhexyl)phthalate	40.000	45.625	-14.1	61	-0.08
84	c	Di-n-octyl phthalate	40.000	45.023	-12.6	65	-0.13
85		Indeno(1,2,3-cd)pyrene	40.000	46.120	-15.3	67	-0.26
86	I	Perylene-d12	20.000	20.000	0.0	59	-0.19
87		Benzo(b)fluoranthene	40.000	52.370	-30.9#	79	-0.16
88		Benzo(k)fluoranthene	40.000	34.329	14.2	52	-0.16
89	C	Benzo(a)pyrene	40.000	44.425	-11.1	67	-0.19
90		Dibenzo(a,h)anthracene	40.000	45.470	-13.7	68	-0.27
91		Benzo(g,h,i)perylene	40.000	45.051	-12.6	69	-0.27

(#) = Out of Range SPCC's out = 0 CCC's out = 1